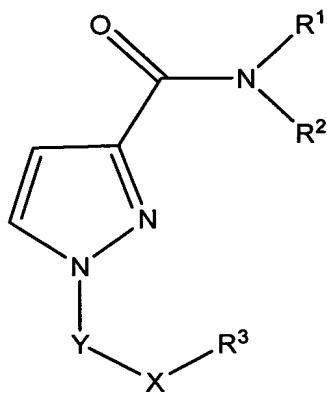


Amendments to the Claims

Please cancel Claims 43 and 45. Please amend Claims 3-34, 36, 38, 40, 42, 44 and 46-52. The Claim Listing below will replace all prior versions of the claims in the application:

Claim Listing

1. (Original) A compound of formula I,



I

wherein either R¹ represents an aryl group or a heteroaryl group, both of which are optionally substituted by one or more substituents selected from G¹ and B¹, which B¹ group may itself be further substituted by one or more substituents selected from G², Z (provided that Z is not directly attached to an aryl or a heteroaryl group) and B² (which B² group is optionally further substituted by one or more substituents selected from G³, B³ and Z, provided that Z is not attached to an aryl or a heteroaryl group); and

R² represents H or C₁₋₆ alkyl, which latter group is optionally substituted by one or more halo groups;

or

when R² represents C₁₋₆ alkyl optionally substituted by halo, R¹ and R² may be linked together forming a further 5- to 7-membered ring, optionally containing 1 to 3 heteroatoms and/or 1 to 3 double bonds, which ring is itself optionally substituted by one or more substituents selected from G¹, Z (provided that the ring is not aromatic in nature) and B¹ (which B¹ group is optionally substituted as described above);

R³ represents C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₈ cycloalkyl, C₃₋₈ heterocycloalkyl, aryl or heteroaryl, all of which groups are optionally substituted by one or more substituents selected from G^{1a}, Z (provided that Z is not directly attached to an aryl or a heteroaryl group) and B¹ (which B¹ group is optionally substituted as described above);

X represents a direct bond, -O- or -N(R⁴)-;

Y represents -C(O)-, -C(S)- or -S(O)₂-;

B¹, B² and B³ independently represent, on each occasion when used above, C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₈ cycloalkyl, C₃₋₈ heterocycloalkyl, aryl or heteroaryl;

G¹, G^{1a}, G² and G³ independently represent, on each occasion when used above, halo, cyano, -N₃, -NO₂, -ONO₂ or -A¹-R⁴;
wherein A¹ represents a spacer group selected from -C(Z)A²-, -N(R⁵)A³-, -OA⁴-, -S- or -S(O)_nA⁵-, in which:

A² represents a single bond, -O-, -S- or -N(R⁵)-;

A³ represents A⁶, -C(Z)N(R⁵)C(Z)N(R⁵)-, -C(Z)N(R⁵)C(Z)O-, -C(Z)N(R⁵)S(O)_nN(R⁵), -C(Z)S-, -S(O)_n-, -S(O)_nN(R⁵)C(Z)N(R⁵)-, -S(O)_nN(R⁵)C(Z)O-, -S(O)_nN(R⁵)S(O)_nN(R⁵)-, -C(Z)O-, -S(O)_nN(R⁵)- or -S(O)_nO-;

A⁴ represents A⁶, -S(O)_n-, -C(Z)O-, -S(O)_nN(R⁵)- or -S(O)_nO-;

A⁵ represents a single bond, -N(R⁵)- or -O-;

A⁶ represents a single bond, -C(Z)- or -C(Z)N(R⁵)-;

Z represents, on each occasion when used above, a substituent connected by a double bond, which is selected from =O, =S, =NR⁴, =NN(R⁴)(R⁵), =NOR⁴, =NS(O)₂N(R⁴)(R⁵), =NCN, =CHNO₂ and =C(R⁴)(R⁵);

R⁴ and R⁵ independently represent, on each occasion when used above, H or B⁴, which B⁴ group is itself optionally substituted by one or more substituents selected from G⁴, Q (provided that Q is not directly attached to an aryl or a heteroaryl group) and B⁵ (which B⁵ group is itself optionally substituted by one or more substituents selected from G⁵, Q (provided that Q is not directly attached to an aryl or a heteroaryl group) and B⁶; or when R⁴ and R⁵ both represent optionally substituted B⁴ groups, then any pair thereof may, for example when present on the same atom or on adjacent atoms, be linked together to form, with those, or other relevant, atoms, a 5- to 7-membered ring, optionally containing 1 to 3 heteroatoms and/or 1 to 3 double bonds, which ring is itself optionally substituted by one or more substituents selected from G⁶, Q (provided that the ring is not aromatic in nature) and B⁴ (which B⁴ group is optionally substituted as described above);

B⁴, B⁵ and B⁶ independently represent on each occasion when used above C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₈ cycloalkyl, C₃₋₈ heterocycloalkyl, aryl or heteroaryl;

G⁴, G⁵ and G⁶ independently represent on each occasion when used above, halo, cyano, N₃, -NO₂, -ONO₂ or -A⁷-R⁶;
 wherein A⁷ represents a spacer group selected from -C(Q)A⁸-, -N(R⁷)A⁹-, -N(R^{7a})A^{9a}-, -OA¹⁰-, -S- or -S(O)_nA¹¹-, in which:
 A⁸ represents a single bond, -O-, -S- or -N(R⁷)-;
 A⁹ represents A¹², -C(Q)S-, -S(O)_n-, -C(Q)O-, -S(O)_nN(R⁷)- or -S(O)_nO-;

A^{9a} represents -C(Q)N(R⁷)C(Q)N(R⁷)-, -C(Q)N(R⁷)C(Q)O-, -C(Q)N(R⁷)S(O)_nN(R⁷)-, -S(O)_nN(R⁷)C(Q)N(R⁷)-, -S(O)_nN(R⁷)C(Q)O-, -S(O)_nN(R⁷)S(O)_nN(R⁷)-;

A¹⁰ represents A¹², -S(O)_n-, -C(Q)O-, -S(O)_nN(R⁷)- or S(O)_nO-;

A¹¹ represents a single bond, -N(R⁷)- or -O-;

A^{12} represents a single bond, $-C(Q)-$ or $-C(Q)N(R^7)-$;

Q represents, on each occasion when used above, a substituent connected by a double bond, which is selected from $=O$, $=S$, $=NR^6$, $=NN(R^6)(R^7)$, $=NOR^6$, $=NS(O)_2N(R^6)(R^7)$, $=NCN$, $=CHNO_2$ and $=C(R^6)(R^7)$;

R^6 , R^7 and R^{7a} independently represent, on each occasion when used above, H , C_{1-6} alkyl, C_{2-6} alkenyl, C_{2-6} alkynyl, C_{3-8} cycloalkyl, C_{3-8} heterocycloalkyl, aryl or heteroaryl, which latter seven groups are optionally substituted by one or more groups selected from halo, C_{1-6} alkyl (optionally substituted by one or more halo groups) - $N(R^8)R^9$, $-OR^8$, $-ONO_2$ and $-SR^8$; or provided that they do not represent H , any pair of R^6 and R^7 may, for example when present on the same atom or on adjacent atoms, be linked together to form, with those, or other relevant, atoms, a 5- to 7-membered ring, optionally containing 1 to 3 heteroatoms and/or 1 to 3 double bonds, which ring is itself optionally substituted by one or more groups selected from halo, C_{1-6} alkyl (optionally substituted by one or more halo groups), $-N(R^8)R^9$, $-OR^8$, $-ONO_2$ and $-SR^8$;

R^8 and R^9 independently represent, on each occasion when used above, H or C_{1-6} alkyl, which latter group is optionally substituted by one or more halo groups; and n represents, on each occasion when used above, 1 or 2;

or a pharmaceutically-acceptable salt thereof,

provided that, when R^2 represents H , Y represents $-C(O)-$ and:

(A) X represents a direct bond and:

- i) R^3 represents phenyl, then R^1 does not represent phenyl, 2-methoxyphenyl, 2-thiazolyl or 6-methyl-2-pyridinyl;
- ii) R^3 represents 4-fluorophenyl, then R^1 does not represent 2-carbomethoxyphenyl, 3-carbomethoxyphenyl or 2,4-dimethylphenyl;
- iii) R^3 represents 2-chlorophenyl, then R^1 does not represent phenyl, 3-bromophenyl or 4-bromophenyl;

- iv) R^3 represents 3-chlorophenyl, then R^1 does not represent phenyl, 2-fluorophenyl, 2-chlorophenyl, 2,3-dichlorophenyl or 2,5-dichlorophenyl;
- v) R^3 represents 4-chlorophenyl, then R^1 does not represent 3-bromophenyl or 4-methoxyphenyl;
- vi) R^3 represents 3-iodophenyl, then R^1 does not represent 2-methoxyphenyl or 2,4-dimethylphenyl;
- vii) R^3 represents 2,4-dichlorophenyl, then R^1 does not represent 4-chlorophenyl or 2,3-dichlorophenyl;
- viii) R^3 represents 3,5-dinitrophenyl, then R^1 does not represent 2,3-dichlorophenyl;
- ix) R^3 represents 2,4-dimethyl-6-oxo-6*H*-pyran-3-yl, then R^1 does not represent 3-carbomethoxyphenyl;
- x) R^3 represents methyl, then R^1 does not represent 3,4-dichlorophenyl, 2-methoxyphenyl, 2-thiazolyl, 4-methyl-2-pyridinyl, 6-methyl-2-pyridinyl or 4-acetylphenyl;
- xi) R^3 represents ethyl, then R^1 does not represent phenyl, 2,3-dichlorophenyl, 4-methoxyphenyl, 2-carbomethoxy-phenyl, 2-thiazolyl or 4-methyl-2-pyridinyl;

(B) X represents -N(H)- and:

- i) R^3 represents phenyl, then R^1 does not represent 4-methoxyphenyl, 2,4-dimethylphenyl or 2-thiazolyl;
- ii) R^3 represents 3-chlorophenyl, then R^1 does not represent 4-methylphenyl;
- iii) R^3 represents 4-chlorophenyl, then R^1 does not represent 3-bromophenyl;
- iv) R^3 represents 3,4-dichlorophenyl, then R^1 does not represent 4-methyl-2-pyridinyl or 6-methyl-2-pyridinyl;
- v) R^3 represents 2'-sulfamoylbiphenyl-4-yl, then R^1 does not represent 5-bromo-2-pyridinyl;
- vi) R^3 represents 1-propyl, then R^1 does not represent phenyl;
- vii) R^3 represents 1-butyl, then R^1 does not represent 4-bromophenyl or 2,4-dimethylphenyl;
- viii) R^3 represents cyclohexyl, then R^1 does not represent 4-methoxyphenyl;

(C) X represents -O- and:

- i) R³ represents phenyl, then R¹ does not represent phenyl or 6-methyl-2-pyridinyl;
- ii) R³ represents methyl, then R¹ does not represent phenyl, 2-fluorophenyl, 2,4-dimethylphenyl, 4-acetylphenyl or 2-thiazolyl;
- iii) R³ represents ethyl, then R¹ does not represent phenyl, 2-fluorophenyl, 4-acetylphenyl or 4-methyl-2-pyridinyl;
- iv) R³ represents 1-butyl, then R¹ does not represent 2-fluorophenyl, 2-methoxyphenyl, 4-methyl-2-pyridinyl or 6-methyl-2-pyridinyl;
- v) R³ represents 2-butyl, then R¹ does not represent 2-thiazolyl or 4-acetylphenyl;
- vi) R³ represents 2-methyl-1-propyl, then R¹ does not represent phenyl or 3-nitrophenyl.

2. (Original) A compound as claimed in Claim 1, wherein R¹ represents an aryl or heteroaryl group, both of which are optionally substituted as defined in Claim 1.
3. (Currently Amended) A compound as claimed in Claim 1 ~~or Claim 2~~, wherein G¹ represents halo, cyano or -A¹-R⁴.
4. (Currently Amended) A compound as claimed in Claim 1 ~~any one of the preceding claims~~, wherein B¹ represents an optionally substituted C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₇ cycloalkyl, C₄₋₇ heterocycloalkyl, or phenyl, group.
5. (Currently Amended) A compound as claimed in Claim 1 ~~any one of the preceding claims~~, wherein G^{1a} represents halo, cyano, -NO₂ or -A¹-R⁴.
6. (Currently Amended) A compound as claimed in Claim 1 ~~any one of the preceding claims~~, wherein G² represents halo, cyano, -ONO₂ or -A¹-R⁴.
7. (Currently Amended) A compound as claimed in Claim 1 ~~in any one of the preceding claims~~, wherein B² represents C₁₋₆ alkyl, C₂₋₆ alkenyl or C₂₋₆ alkynyl, all of which are optionally substituted by one or more G³ and/or B³ groups.

8. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein G³ represents halo, -ONO₂, -N(R⁵)(R⁴) or -OR⁴.
9. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein B³ represents C₁₋₆ alkyl, C₂₋₆ alkenyl or C₂₋₆ alkynyl.
10. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein when A¹ represents -N(R⁵)A³-, A³ represents A⁶, -C(Z)S-, -S(O)_n-, -C(Z)O- or -S(O)_nN(R⁵)-.
11. (Currently Amended) A compound as claimed in any one of Claims Claim 1 to 9, wherein when A¹ represents -OA⁴-, A⁴ represents A⁶.
12. (Currently Amended) A compound as claimed in any one of Claims Claim 1 to 9, wherein when A¹ represents -S(O)_nA⁵-, A⁵ represents a single bond or -N(R⁵)-.
13. (Currently Amended) A compound as claimed in any one of Claims Claim 1 to 9, wherein when A¹ represents -C(Z)A²-, A² represents a single bond, -O- or -N(R⁵)-.
14. (Currently Amended) A compound as claimed in any one of Claims Claim 1 to 11 or 13 wherein A¹ represents -C(Z)A²-, -N(R⁵)A³- or -OA⁴-.
15. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein Z represents =O or =NR⁴.
16. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein when any pair of R⁴ and R⁵ are linked together to form a ring, they are optionally substituted with G⁶ and/or B⁴.
17. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein G⁴ represents halo, cyano, -ONO₂ or -A⁷-R⁶.

18. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein B⁵ represents C₁₋₆ alkyl, C₂₋₆ alkenyl or C₂₋₆ alkynyl, all of which are optionally substituted by one or more G⁵ and/or B⁶ groups.
19. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein G⁵ represents halo, -ONO₂, -N(R⁷)(R⁶) or -OR⁶.
20. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein B⁶ represents C₁₋₆ alkyl, C₂₋₆ alkenyl or C₂₋₆ alkynyl.
21. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein G⁶ represents halo, cyano or -A⁷ -R⁶.
22. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein A⁷ represents -C(Q)A⁸-, -N(R⁷)A⁹-, -OA¹⁰-, -S- or -S(O)_nA¹¹-.
23. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein when A⁷ represents -N(R⁷)A⁹-, A⁹ represents A¹², -C(Q)S-, -S(O)_n-, -C(Q)O- or -S(O)_nN(R⁷)-.
24. (Currently Amended) A compound as claimed in Claim 1 any one of Claims 1 to 22, wherein when A⁷ represents -OA¹⁰-, A¹⁰ represents A¹².
25. (Currently Amended) A compound as claimed in Claim 1 any one of Claims 1 to 22, wherein when A⁷ represents -S(O)_nA¹¹-, A¹¹ represents a single bond or -N(R⁷)-.
26. (Currently Amended) A compound as claimed in Claim 1 any one of Claims 1 to 22, wherein when A⁷ represents -C(Q)A⁸-, A⁸ represents a single bond, -O- or -N(R⁷)-.
27. (Currently Amended) A compound as claimed in Claim 1 any one of the preceding claims, wherein Q represents =O or =NR⁶.

28. (Currently Amended) A compound as claimed in Claim 1 ~~any one of the preceding claims~~, wherein R⁶, R⁷ and R^{7a} independently represent H, C₁₋₆ alkyl, C₂₋₆ alkenyl or C₂₋₆ alkynyl, all of which groups are optionally substituted by one or more groups selected from halo, C₁₋₆ alkyl, -N(R⁸)R⁹, OR⁸ and -ONO₂.
29. (Currently Amended) A compound as claimed in Claim 1 ~~any one of Claims 1 to 27~~ wherein when any pair of R⁶ and R⁷ are linked together to form a ring, that ring is optionally substituted by one or more groups selected from halo, C₁₋₆ alkyl (optionally substituted by one or more halo groups), -N(R⁸)R⁹, -OR⁸ and -ONO₂.
30. (Currently Amended) A compound as claimed in Claim 1 ~~any one of the preceding claims~~, wherein B⁴ represents an optionally substituted C₁₋₆ alkyl, C₂₋₆ alkenyl, C₂₋₆ alkynyl, C₃₋₇ cycloalkyl, C₄₋₇ heterocycloalkyl, or phenyl, group.
31. (Currently Amended) A compound as claimed in Claim 1 ~~any one of the preceding claims~~ wherein R⁴ and/or R⁵ independently represent H or C₁₋₆ alkyl, which latter group is optionally substituted by one or more fluoro groups.
32. (Currently Amended) A compound as claimed in Claim 1 ~~any one of the preceding claims~~, wherein X represents a direct bond, -O-, -N(H)- or -N(Me)-.
33. (Currently Amended) A compound as claimed in Claim 1 ~~any one of the preceding claims~~ wherein R² represents H, methyl or ethyl.
34. (Currently Amended) A compound as claimed in Claim 1 ~~any one of Claims 1, 32 or 33~~, wherein R¹ represents an optionally substituted phenyl, naphthyl, pyrrolidinyl, piperidinyl, pyrrolyl, furanyl, thiophenyl, pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, thiazolyl, pyridinyl, indazolyl, indolyl, indolinyl, isoindolinyl, oxindolyl, quinolinyl, 1,2,3,4-tetrahydroquinolinyl, isoquinolinyl, 1,2,3,4-tetrahydroisoquinolinyl, quinolizinyl, benzofuranyl, isobenzofuranyl, chromanyl, benzothiophenyl, pyridazinyl, pyrimidinyl, pyrazinyl, indazolyl, benzimidazolyl, quinazolinyl, quinoxalinyl, 1,3-benzodioxolyl, benzothiazolyl, or benzodioxanyl, group.

35. (Original) A compound as claimed in Claim 34, wherein R¹ represents optionally substituted phenyl, 2-pyridinyl, 3-pyridinyl, 2-thiophenyl, 4-pyrazolyl, 5-isoxazolyl, 1,3-benzodioxolyl, indazolyl, benzothiazolyl, or quinolinyl, group.
36. (Currently Amended) A compound as claimed in Claim 34 or ~~Claim 35~~, wherein the optional substituent(s) are selected from halo, cyano, C₁₋₆ alkyl (which alkyl group may be linear or branched, and/or substituted by one or more fluoro and/or C₃₋₆ cycloalkyl groups), C₂₋₆ alkenyl, C₃₋₆ cycloalkyl, phenyl, pyrrolidinyl, piperidinyl, piperazinyl, tetrahydrofuranyl, tetrahydropyranyl, morpholinyl, thiomethyl, methylsulfinyl, methylsulfonyl, -OR¹⁰, -N(R¹⁰)R¹¹, -C(O)OR¹⁰, -C(O)R¹⁰, -C(O)N(R¹⁰)R¹¹, -S(O)₂N(R¹⁰)R¹¹ and -N(R¹⁰)S(O)₂R¹², wherein R¹⁰ and R¹¹ independently represent H, phenyl, C₁₋₆ alkyl (which alkyl group is optionally substituted by one or more fluoro atom), C₂₋₆ alkenyl or C₃₋₆ cycloalkyl; or R¹⁰ and R¹¹ may be linked together to form, with the nitrogen atom to which they are attached, a 5- to 7-membered ring, optionally containing one additional heteroatom and optionally substituted with one or more C₁₋₆ alkyl groups, which alkyl groups are themselves optionally substituted by one or more halo groups; and R¹² represents phenyl, C₁₋₆ alkyl (which alkyl group is optionally substituted by one or more fluoro atom), C₂₋₆ alkenyl or C₃₋₆ cycloalkyl.
37. (Original) A compound as claimed in Claim 36, wherein the optional substituent(s) are selected from carbomethoxy, methyl, dimethylamino, cyano, cholor, fluoro, trifluoromethyl, bromo, methoxy and trifluoromethoxy.
38. (Currently Amended) A compound as claimed in Claim 1 ~~any one of Claims 1 or 32 to 37~~, wherein R³ represents an optionally substituted C₁₋₆ alkyl, C₃₋₆ cycloalkyl, phenyl, naphthyl, pyrrolidinyl, piperidinyl, piperazinyl, pyrrolyl, furanyl, thiophenyl, pyrazolyl, imidazolyl, oxazolyl, isoxazolyl, thiazolyl, pyridinyl, indazolyl, indolyl, indolinyl, isoindolinyl, oxindolyl, quinolinyl, 1,2,3,4-tetrahydroquinolinyl, isoquinolinyl, 1,2,3,4-tetrahydroisoquinolinyl, quinolizinyl, benzofuranyl, isobenzofuranyl, chromanyl, benzothiophenyl, pyridazinyl, pyrimidinyl, pyrazinyl,

indazolyl, benzimidazolyl, quinazolinyl, quinoxalinyl, 1,3-benzodioxolyl, benzothiazolyl, or benzodioxanyl, group.

39. (Original) A compound as claimed in Claim 38, wherein R³ represents an optionally substituted C₁₋₆ alkyl, cyclohexyl, phenyl, 2-thiophenyl, 2-furanyl, 3-furanyl, 2-pyrrolyl, 1-naphthyl, 4-piperazinyl, 4-piperidinyl, benzofuranyl, or 1,3-benzodioxolyl, group.
40. (Currently Amended) A compound as claimed in Claim 38 or ~~Claim 39~~, wherein the optional substituent(s) are selected from halo, -NO₂, cyano, C₁₋₆ alkyl (which alkyl group may be linear or branched, and/or optionally substituted with one or more halo, C₁₋₆ alkyl, C₂₋₆ alkenyl and/or C₃₋₆ cycloalkyl, groups, which latter three groups are themselves optionally substituted with one or more halo and/or C₁₋₆ alkyl groups), C₂₋₆ alkenyl (optionally substituted with one or more C₁₋₆ alkyl groups), C₃₋₆ cycloalkyl (optionally substituted with one or more halo groups), phenyl (optionally substituted with one or more halo groups), pyrrolidinyl, piperidinyl, piperazinyl, tetrahydrofuranyl, tetrahydropyranyl, morpholinyl, thiomethyl, methylsufinyl, methylsulfonyl, =O, -OR¹³, -N(R¹³)R¹⁴, -C(O)OR¹³, -C(O)R¹³, -C(O)N(R¹³)R¹⁴, -S(O)₂N(R¹³)R¹⁴ and -N(R¹³)S(O)₂R¹⁵, wherein R¹³ and R¹⁴ independently represent H, phenyl, C₁₋₆ alkyl (which alkyl group is optionally substituted by one or more fluoro atom), C₂₋₆ alkenyl or C₃₋₆ cycloalkyl; or R¹³ and R¹⁴ may be linked together to form, with the nitrogen atom to which they are attached, a 5- to 7- membered ring, optionally containing one additional heteroatom and optionally substituted with one or more C₁₋₆ alkyl groups, which alkyl groups are themselves optionally substituted by one or more halo groups; and R¹⁵ represents phenyl, C₁₋₆ alkyl (which alkyl group is optionally substituted by one or more fluoro atom), C₂₋₆ alkenyl or C₃₋₆ cycloalkyl.
41. (Original) A compound as claimed in Claim 40, wherein the optional substituent(s) are selected from methyl, ethyl, ethoxy, trifluoromethyl, fluoro, chloro, iodo, phenyl, 2-chlorophenyl, 4-chlorophenyl, *n*-pentyl, *i*-propyl, nitro, *t*-butyl, -CH₂CH=CHC₈H₁₇, trifluoroacetyl, carbomethoxy, carboethoxy and trifluoromethoxy.

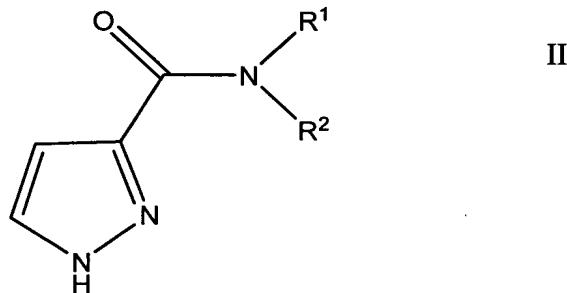
42. (Currently Amended) A compound as claimed in Claim 1 any one of Claims 1 to 32 to 41, wherein R¹ is phenyl, 2-chlorophenyl, 2-chloro-4-fluorophenyl, 3-chloro-4-fluorophenyl, 2,6-dichlorophenyl, 5-chloro-2-cyanophenyl, 2-fluoro-5-trifluoromethylphenyl, 2-bromo-4-trifluoromethoxyphenyl, 2-methoxy-6-methylphenyl, 3-cyanophenyl, 4-trifluoromethylphenyl, 4-dimethylaminophenyl, 4-carbomethoxyphenyl, 1,3,5-trimethyl-1*H*-pyrazol-4-yl, 3-methylisoxazol-5-yl, 3-pyridinyl, 2-chloro-3-pyridinyl, 3-methyl-2-pyridinyl, 3-carbomethoxythiophen-2-yl or 1,3-benzodioxolyl;
R² is hydrogen or methyl;
R³ is methyl, *n*-butyl, *n*-pentyl, 1-octyl, oleoyl, (1*R*,2*S*,5*R*)-(-)-menthyl, 2-chlorobenzyl, benzyl, phenyl, 3-fluorophenyl, 3-chlorophenyl, 4-chlorophenyl, 2-fluoro-5-iodophenyl, 5-fluoro-2-methylphenyl, 4-*tert*-butyl-phenyl, 4-pentylphenyl, 3-trifluoromethylphenyl, 4-trifluoromethoxyphenyl, 4-nitrophenyl, 2-ethoxyphenyl, 1-naphthyl, 2-furanyl, 2,5-dimethyl-3-furanyl, 2-carbomethoxy-5-furanyl, 1-methyl-1*H*-pyrrol-2-yl, 3-methyl-2-benzofuranyl, 3-methyl-2-thiophenyl, 1(*N*)-methyl-4-piperazinyl, 1(*N*)-(2,2,2-trifluoroacetyl)piperidin-4-yl, ethylhexanoate or 1,3-benzodioxolyl; Y is -C(O)-, -C(S)- or -S(O)₂-; and
X is a bond, -N(H)-, -N(Me)-, or -O-.
43. (Cancelled)
44. (Currently Amended) A pharmaceutical formulation including a compound of formula I, as defined in Claim 1 any one of Claims 1 to 42, but without the provisos, or a pharmaceutically acceptable salt thereof, in admixture with a pharmaceutically acceptable adjuvant, diluent or carrier.
45. (Cancelled)
46. (Currently Amended) A use The method as claimed in Claim 45 49 wherein the lipoxygenase is 15-lipoxygenase.

47. (Currently Amended) ~~A use~~ The method as claimed in ~~Claim 45~~ or Claim 46, wherein the disease is inflammation and/or has an inflammatory component.
48. (Currently Amended) The method ~~A use~~ as claimed in Claim 47 wherein the inflammatory disease is asthma, chronic obstructive pulmonary disease (COPD), pulmonary fibrosis, an allergic disorder, rhinitis, inflammatory bowel disease, an ulcer, inflammatory pain, fever, atherosclerosis, coronary artery disease, vasculitis, pancreatitis, arthritis, osteoarthritis, rheumatoid arthritis, conjunctivitis, iritis, scleritis, uveitis, a wound, dermatitis, eczema, psoriasis, stroke, diabetes, autoimmune disease, Alzheimer's disease, multiple sclerosis, sarcoidosis, Hodgkin's disease or another malignancy.
49. (Currently Amended) A method of treatment of a disease in which inhibition of the activity of a lipoxygenase is desired and/or required, which method comprises administration of a therapeutically effective amount of a compound of formula I as defined in Claim 1 ~~any one of Claims 1 to 42~~, but without the provisos, or a pharmaceutically-acceptable salt thereof, to a patient suffering from, or susceptible to, such a condition.
50. (Currently Amended) A combination product comprising:
 - (A) a compound of formula I as defined in Claim 1 ~~any one of Claims 1 to 42~~, but without the provisos; and
 - (B) another therapeutic agent that is useful in the treatment of inflammation, wherein each of components (A) and (B) is formulated in admixture with a pharmaceutically-acceptable adjuvant, diluent or carrier.
51. (Currently Amended) A combination product as claimed in Claim 50 which comprises a pharmaceutical formulation including ~~a~~ the compound of formula I ~~as defined in any one of Claims 1 to 42, but without the provisos~~, another therapeutic agent that is useful in the treatment of inflammation, and a pharmaceutically-acceptable adjuvant, diluent or carrier.

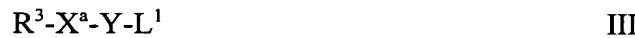
52. (Currently Amended) A combination product as claimed in Claim 50 which comprises a kit of parts comprising components:
- (a) a pharmaceutical formulation including a the compound of formula I as defined in any one of Claims 1 to 42, but without the provisos, in admixture with a pharmaceutically-acceptable adjuvant, diluent or carrier; and
 - (b) a pharmaceutical formulation including another therapeutic agent that is useful in the treatment of inflammation in admixture with a pharmaceutically-acceptable adjuvant, diluent or carrier,
- which components (a) and (b) are each provided in a form that is suitable for administration in conjunction with the other.

53. (Original) A process for the preparation of a compound as defined in Claim 1, which comprises:

- (i) for compounds of formula I in which, when Y is $-S(O)_2-$, X represents a direct bond or $-N(R^4)-$, in which R^4 represents B^4 , reaction of a compound of formula II,

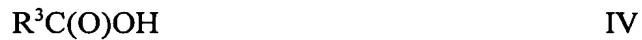


wherein R^1 and R^2 are as defined in Claim 1, with a compound of formula III,



wherein X^8 represents a direct bond or $-N(B^4)-$ when Y represents $-S(O)_2-$ or, for all other values of Y , represents X as defined in Claim 1, R^3 and Y are as defined in Claim 1 and L^1 represents a suitable leaving group;

(ii) for compounds of formula I in which X represents a single bond and Y represents $-C(O)-$, reaction of a compound of formula II as defined above with a compound of formula IV,



wherein R^3 is as defined in Claim 1;

(iii) for compounds of formula I in which X represents a direct bond and Y represents a $-C(O)-$ or a $-C(S)-$ group, reaction of a compound of formula II as defined above with a compound of formula V,



wherein Y^8 represents $-C(O)-$ or $-C(S)-$ and R^3 is as defined in Claim 1;

(iv) for compounds of formula I, in which X represents $-NH-$ and Y represents $-C(O)-$ or $-C(S)-$, reaction of a compound of formula II as defined above with a compound of formula VI,



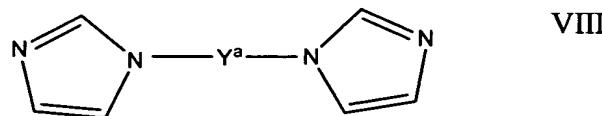
wherein R^3 is as defined in Claim 1 and Y^a is as defined above;

(v) for compounds of formula I in which Y represents $-C(O)-$ or $-C(S)-$, reaction of a compound of formula II as defined above with:

(a) a compound of formula VII,



(b) a compound of formula VIII,



wherein, in both cases, Y^a is as defined above; or

(c) when Y represents $-C(O)-$, triphosgene, followed by:

(1) for compounds of formula I in which X represents a direct bond, reaction with a compound of formula IX,



wherein M represents a metal such as Mn, Fe, Ni, Cu, Zn, Pd or Ce, or a salt or complex thereof and R^3 is as defined in Claim 1;

(2) for compounds of formula I wherein X represents O, reaction with a compound of formula X,



wherein R^3 is as defined in Claim 1; or

(3) for compounds of formula I wherein X represents $-N(R^4)-$, reaction with a compound of formula XI,



wherein R^3 and R^4 are as defined in Claim 1;

(vi) for compounds of formula I in which X represents $-N(R^4)-$ and R^4 is other than hydrogen, reaction of a corresponding compound of formula I in which X represents $-N(H)-$ with a compound of formula XII,

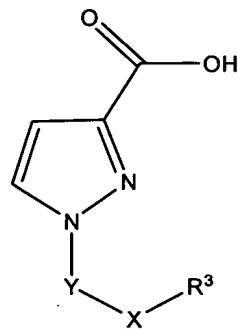
$R^4 \cdot L^1$

XII

wherein R^4 is as defined in Claim 1 and L^1 is as defined above;

(vii) for compounds of formula I in which Y represents $-C(S)-$, reaction of a corresponding compound of formula I in which Y represents $-C(O)-$ with a suitable reagent for the conversion of a carbonyl group to a thiocarbonyl group;

(viii) reaction of a compound of formula XIII,



XIII

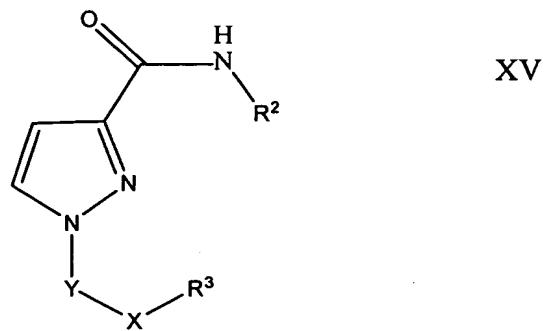
wherein R^3 , Y and X are defined in Claim 1, with a compound of formula XIV,

 $HN(R^1)(R^2)$

XIV

wherein R^1 and R^2 are as defined in Claim 1; or

(ix) reaction of a compound of formula XV,



XV

wherein R², R³, Y and X are as defined in Claim 1, with a compound of formula XVI,



XVI

wherein L² represents a suitable leaving group and R¹ is as defined in Claim 1.